

- ☒ Final Report  
☐ Re-Issued Report  
☐ Revised Report

Report Date:  
10-Nov-16 15:07

## Laboratory Report

Gulf Oil L.P.  
281 Eastern Avenue  
Chelsea, MA 02150  
Attn: Andrew P. Adams

Project: Gulf Terminal - Chelsea, MA  
Project #: Gulf Chelsea

<u>Laboratory ID</u>	<u>Client Sample ID</u>	<u>Matrix</u>	<u>Date Sampled</u>	<u>Date Received</u>
SC27780-01	Outfall 003	Surface Water	28-Oct-16 12:30	31-Oct-16 16:44

I attest that the information contained within the report has been reviewed for accuracy and checked against the quality control requirements for each method. These results relate only to the sample(s) as received.  
All applicable NELAC requirements have been met.

Massachusetts # M-MA138/MA1110  
Connecticut # PH-0777  
Florida # E87936  
Maine # MA138  
New Hampshire # 2972/2538  
New Jersey # MA011  
New York # 11393  
Pennsylvania # 68-04426/68-02924  
Rhode Island # LAO00348  
USDA # P330-15-00375  
Vermont # VT-11393



Authorized by:



June O'Connor  
Laboratory Director

Eurofins Spectrum Analytical holds primary certification in the State of Massachusetts for the analytes as indicated with an X in the "Cert." column within this report. Please note that the State of Massachusetts does not offer certification for all analytes. Please refer to our website for specific certification holdings in each state.

Please note that this report contains 11 pages of analytical data plus Chain of Custody document(s). When the Laboratory Report is indicated as revised, this report supersedes any previously dated reports for the laboratory ID(s) referenced above. Where this report identifies subcontracted analyses, copies of the subcontractor's test report are available upon request. This report may not be reproduced, except in full, without written approval from Eurofins Spectrum Analytical, Inc.

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*Please contact the Laboratory or Technical Director at 800-789-9115 with any questions regarding the data contained in this laboratory report.*

**CASE NARRATIVE:**

Data has been reported to the RDL. This report excludes estimated concentrations detected below the RDL and above the MDL (J-Flag).

All non-detects and all results below the reporting limit are reported as "<" (less than) the reporting limit in this report.

The samples were received 0.8 degrees Celsius, please refer to the Chain of Custody for details specific to temperature upon receipt. An infrared thermometer with a tolerance of +/- 1.0 degrees Celsius was used immediately upon receipt of the samples.

If a Matrix Spike (MS), Matrix Spike Duplicate (MSD) or Duplicate (DUP) was not requested on the Chain of Custody, method criteria may have been fulfilled with a source sample not of this Sample Delivery Group.

Analyses for Total Hardness, pH, and Total Residual Chlorine fall under the state of Pennsylvania code Chapter 252.6 accreditation by rule.

**See below for any non-conformances and issues relating to quality control samples and/or sample analysis/matrix.**

**SW846 8260C****Calibration:**

1611006

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Analyte quantified by quadratic equation type calibration.

Naphthalene

This affected the following samples:

1619121-BLK1

1619121-BS1

1619121-BSD1

Outfall 003

S609436-CCV1

S609464-ICV1

## Sample Acceptance Check Form

Client: Gulf Oil L.P.  
Project: Gulf Terminal - Chelsea, MA / Gulf Chelsea  
Work Order: SC27780  
Sample(s) received on: 10/31/2016

*The following outlines the condition of samples for the attached Chain of Custody upon receipt.*

	<u>Yes</u>	<u>No</u>	<u>N/A</u>
Were custody seals present?	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>
Were custody seals intact?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
Were samples received at a temperature of $\leq 6^{\circ}\text{C}$ ?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Were samples refrigerated upon transfer to laboratory representative?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Were sample containers received intact?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Were samples properly labeled (labels affixed to sample containers and include sample ID, site location, and/or project number and the collection date)?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Were samples accompanied by a Chain of Custody document?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Does Chain of Custody document include proper, full, and complete documentation, which shall include sample ID, site location, and/or project number, date and time of collection, collector's name, preservation type, sample matrix and any special remarks concerning the sample?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Did sample container labels agree with Chain of Custody document?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Were samples received within method-specific holding times?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>

## Summary of Hits

**Lab ID:** SC27780-01

**Client ID:** Outfall 003

Parameter	Result	Flag	Reporting Limit	Units	Analytical Method
Oil and Grease by EPA 1664A	< 1.4		1.4	mg/L	E1664A
Total Suspended Solids	1.9		0.5	mg/l	SM2540D (11)

*Please note that because there are no reporting limits associated with hazardous waste characterizations or micro analyses, this summary does not include hits from these analyses if included in this work order.*

Sample Identification

Outfall 003

SC27780-01

Client Project #

Gulf Chelsea

Matrix

Surface Water

Collection Date/Time

28-Oct-16 12:30

Received

31-Oct-16

<i>CAS No.</i>	<i>Analyte(s)</i>	<i>Result</i>	<i>Flag</i>	<i>Units</i>	<i>*RDL</i>	<i>MDL</i>	<i>Dilution</i>	<i>Method Ref.</i>	<i>Prepared</i>	<i>Analyzed</i>	<i>Analyst</i>	<i>Batch</i>	<i>Cert.</i>
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**Volatile Organic Compounds**

Volatile Organic Aromatics by SW846 8260

Prepared by method SW846 5030 Water MS

71-43-2	Benzene	< 1.0		µg/l	1.0	0.3	1	SW846 8260C	02-Nov-16	02-Nov-16	TS	1619121	
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91-20-3	Naphthalene	< 1.0		µg/l	1.0	0.3	1	"	"	"	"	"	
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*Surrogate recoveries:*

460-00-4	4-Bromofluorobenzene	99			70-130 %			"	"	"	"	"	
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2037-26-5	Toluene-d8	99			70-130 %			"	"	"	"	"	
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17060-07-0	1,2-Dichloroethane-d4	102			70-130 %			"	"	"	"	"	
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1868-53-7	Dibromofluoromethane	97			70-130 %			"	"	"	"	"	
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**Semivolatile Organic Compounds by GCMS**SVOCs by SIMPrepared by method SW846 3510C

50-32-8	Benzo (a) pyrene	< 0.051		µg/l	0.051	0.037	1	SW846 8270D SIM	03-Nov-16	07-Nov-16	MSL	1619194	
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91-20-3	Naphthalene	< 0.051		µg/l	0.051	0.027	1	"	"	"	"	"	
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*Surrogate recoveries:*

205440-82-0	Benzo (e) pyrene-d12	50			30-130 %			"	"	"	"	"	
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**General Chemistry Parameters**

pH	7.60		pH Units				1	ASTM D 1293-99B	01-Nov-16 10:40	01-Nov-16 10:40	TY	1619080	X
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Total Suspended Solids	1.9		mg/l	0.5	0.2		1	SM2540D (11)	02-Nov-16	03-Nov-16	CMB	1619128	X
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**Subcontracted Analyses**Prepared by method 365564*Analysis performed by Phoenix Environmental Labs, Inc. \* - MACT007*

Oil and Grease by EPA 1664A	< 1.4		mg/L	1.4	1.4		1	E1664A		04-Nov-16 08:19	MACT0	365564A	
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# Volatile Organic Compounds - Quality Control

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
<b>Batch 1619121 - SW846 5030 Water MS</b>										
<b><u>Blank (1619121-BLK1)</u></b>					<u>Prepared &amp; Analyzed: 02-Nov-16</u>					
Benzene	< 1.0		µg/l	1.0						
Ethylbenzene	< 1.0		µg/l	1.0						
Methyl tert-butyl ether	< 1.0		µg/l	1.0						
Naphthalene	< 1.0		µg/l	1.0						
Toluene	< 1.0		µg/l	1.0						
m,p-Xylene	< 2.0		µg/l	2.0						
o-Xylene	< 1.0		µg/l	1.0						
Surrogate: 4-Bromofluorobenzene	49.2		µg/l		50.0		98	70-130		
Surrogate: Toluene-d8	50.3		µg/l		50.0		101	70-130		
Surrogate: 1,2-Dichloroethane-d4	51.3		µg/l		50.0		103	70-130		
Surrogate: Dibromofluoromethane	47.8		µg/l		50.0		96	70-130		
<b><u>LCS (1619121-BS1)</u></b>					<u>Prepared &amp; Analyzed: 02-Nov-16</u>					
Benzene	20.9		µg/l		20.0		105	70-130		
Ethylbenzene	21.4		µg/l		20.0		107	70-130		
Methyl tert-butyl ether	19.8		µg/l		20.0		99	70-130		
Naphthalene	20.7		µg/l		20.0		103	70-130		
Toluene	20.0		µg/l		20.0		100	70-130		
m,p-Xylene	21.2		µg/l		20.0		106	70-130		
o-Xylene	21.5		µg/l		20.0		107	70-130		
Surrogate: 4-Bromofluorobenzene	51.6		µg/l		50.0		103	70-130		
Surrogate: Toluene-d8	49.8		µg/l		50.0		100	70-130		
Surrogate: 1,2-Dichloroethane-d4	50.8		µg/l		50.0		102	70-130		
Surrogate: Dibromofluoromethane	51.6		µg/l		50.0		103	70-130		
<b><u>LCS Dup (1619121-BSD1)</u></b>					<u>Prepared &amp; Analyzed: 02-Nov-16</u>					
Benzene	20.5		µg/l		20.0		102	70-130	2	20
Ethylbenzene	21.2		µg/l		20.0		106	70-130	0.8	20
Methyl tert-butyl ether	19.6		µg/l		20.0		98	70-130	0.9	20
Naphthalene	21.0		µg/l		20.0		105	70-130	2	20
Toluene	20.3		µg/l		20.0		102	70-130	2	20
m,p-Xylene	21.4		µg/l		20.0		107	70-130	1	20
o-Xylene	21.4		µg/l		20.0		107	70-130	0.4	20
Surrogate: 4-Bromofluorobenzene	51.6		µg/l		50.0		103	70-130		
Surrogate: Toluene-d8	49.9		µg/l		50.0		100	70-130		
Surrogate: 1,2-Dichloroethane-d4	49.8		µg/l		50.0		100	70-130		
Surrogate: Dibromofluoromethane	51.6		µg/l		50.0		103	70-130		

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## Semivolatile Organic Compounds by GCMS - Quality Control

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
<b>Batch 1619194 - SW846 3510C</b>										
<b><u>Blank (1619194-BLK2)</u></b>					<u>Prepared: 03-Nov-16 Analyzed: 04-Nov-16</u>					
Acenaphthene	< 0.050		µg/l	0.050						
Acenaphthylene	< 0.050		µg/l	0.050						
Anthracene	< 0.050		µg/l	0.050						
Benzo (a) anthracene	< 0.050		µg/l	0.050						
Benzo (a) pyrene	< 0.050		µg/l	0.050						
Benzo (b) fluoranthene	< 0.050		µg/l	0.050						
Benzo (g,h,i) perylene	< 0.050		µg/l	0.050						
Benzo (k) fluoranthene	< 0.050		µg/l	0.050						
Chrysene	< 0.050		µg/l	0.050						
Dibenzo (a,h) anthracene	< 0.050		µg/l	0.050						
Fluoranthene	< 0.050		µg/l	0.050						
Fluorene	< 0.050		µg/l	0.050						
Indeno (1,2,3-cd) pyrene	< 0.050		µg/l	0.050						
Naphthalene	< 0.050		µg/l	0.050						
Phenanthrene	< 0.050		µg/l	0.050						
Pyrene	< 0.050		µg/l	0.050						
<i>Surrogate: Benzo (e) pyrene-d12</i>	<i>0.700</i>		µg/l		<i>1.00</i>		<i>70</i>	<i>30-130</i>		
<b><u>LCS (1619194-BS2)</u></b>					<u>Prepared: 03-Nov-16 Analyzed: 04-Nov-16</u>					
Acenaphthene	<b>0.764</b>		µg/l	0.050	1.00		76	40-140		
Acenaphthylene	<b>0.725</b>		µg/l	0.050	1.00		72	40-140		
Anthracene	<b>0.833</b>		µg/l	0.050	1.00		83	40-140		
Benzo (a) anthracene	<b>0.834</b>		µg/l	0.050	1.00		83	40-140		
Benzo (a) pyrene	<b>0.864</b>		µg/l	0.050	1.00		86	40-140		
Benzo (b) fluoranthene	<b>0.724</b>		µg/l	0.050	1.00		72	40-140		
Benzo (g,h,i) perylene	<b>0.800</b>		µg/l	0.050	1.00		80	40-140		
Benzo (k) fluoranthene	<b>0.883</b>		µg/l	0.050	1.00		88	40-140		
Chrysene	<b>0.889</b>		µg/l	0.050	1.00		89	40-140		
Dibenzo (a,h) anthracene	<b>0.838</b>		µg/l	0.050	1.00		84	40-140		
Fluoranthene	<b>0.839</b>		µg/l	0.050	1.00		84	40-140		
Fluorene	<b>0.776</b>		µg/l	0.050	1.00		78	40-140		
Indeno (1,2,3-cd) pyrene	<b>0.875</b>		µg/l	0.050	1.00		88	40-140		
Naphthalene	<b>0.743</b>		µg/l	0.050	1.00		74	40-140		
Phenanthrene	<b>0.754</b>		µg/l	0.050	1.00		75	40-140		
Pyrene	<b>0.847</b>		µg/l	0.050	1.00		85	40-140		
<i>Surrogate: Benzo (e) pyrene-d12</i>	<i>0.700</i>		µg/l		<i>1.00</i>		<i>70</i>	<i>30-130</i>		
<b><u>LCS Dup (1619194-BS2)</u></b>					<u>Prepared: 03-Nov-16 Analyzed: 04-Nov-16</u>					
Acenaphthene	<b>0.661</b>		µg/l	0.050	1.00		66	40-140	14	20
Acenaphthylene	<b>0.664</b>		µg/l	0.050	1.00		66	40-140	9	20
Anthracene	<b>0.754</b>		µg/l	0.050	1.00		75	40-140	10	20
Benzo (a) anthracene	<b>0.808</b>		µg/l	0.050	1.00		81	40-140	3	20
Benzo (a) pyrene	<b>0.742</b>		µg/l	0.050	1.00		74	40-140	15	20
Benzo (b) fluoranthene	<b>0.747</b>		µg/l	0.050	1.00		75	40-140	3	20
Benzo (g,h,i) perylene	<b>0.647</b>	QR2	µg/l	0.050	1.00		65	40-140	21	20
Benzo (k) fluoranthene	<b>0.743</b>		µg/l	0.050	1.00		74	40-140	17	20
Chrysene	<b>0.781</b>		µg/l	0.050	1.00		78	40-140	13	20
Dibenzo (a,h) anthracene	<b>0.708</b>		µg/l	0.050	1.00		71	40-140	17	20
Fluoranthene	<b>0.761</b>		µg/l	0.050	1.00		76	40-140	10	20
Fluorene	<b>0.664</b>		µg/l	0.050	1.00		66	40-140	16	20
Indeno (1,2,3-cd) pyrene	<b>0.760</b>		µg/l	0.050	1.00		76	40-140	14	20
Naphthalene	<b>0.702</b>		µg/l	0.050	1.00		70	40-140	6	20
Phenanthrene	<b>0.694</b>		µg/l	0.050	1.00		69	40-140	8	20

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# Semivolatile Organic Compounds by GCMS - Quality Control

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
<b>Batch 1619194 - SW846 3510C</b>										
<b><u>LCS Dup (1619194-BSD2)</u></b>					<u>Prepared: 03-Nov-16 Analyzed: 04-Nov-16</u>					
Pyrene	<b>0.767</b>		µg/l	0.050	1.00		77	40-140	10	20
Surrogate: Benzo (e) pyrene-d12	0.630		µg/l		1.00		63	30-130		



# General Chemistry Parameters - Quality Control

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
<b>Batch 1619080 - General Preparation</b>										
<b><u>Duplicate (1619080-DUP1)</u></b>										
pH	7.56		pH Units			7.60			0.5	5
<b><u>Reference (1619080-SRM1)</u></b>										
pH	6.05		pH Units		6.00		101	97.5-102.5		
<b><u>Reference (1619080-SRM2)</u></b>										
pH	6.04		pH Units		6.00		101	97.5-102.5		
<b>Batch 1619128 - General Preparation</b>										
<b><u>Blank (1619128-BLK1)</u></b>										
Total Suspended Solids	< 0.5		mg/l	0.5						
<b><u>LCS (1619128-BS1)</u></b>										
Total Suspended Solids	102		mg/l	10.0	100		102	90-110		

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## Subcontracted Analyses - Quality Control

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
<b>Batch 365564A - 365564</b>										
<b><u>BLK (BV72754-BLK)</u></b>					<u>Prepared &amp; Analyzed: 04-Nov-16</u>					
Oil and Grease by EPA 1664A	< 1.4		mg/L	1.4	40			-		
<b><u>LCS (BV72754-LCS)</u></b>					<u>Prepared: Analyzed: 04-Nov-16</u>					
Oil and Grease by EPA 1664A	<b>39.10</b>		mg/L	1.4	40		98	85-115		20

## Notes and Definitions

QR2	The RPD result exceeded the QC control limits; however, both percent recoveries were acceptable. Sample results for the QC batch were accepted based on percent recoveries and completeness of QC data.
dry	Sample results reported on a dry weight basis
NR	Not Reported
RPD	Relative Percent Difference
OG	The required Matrix Spike and Matrix Spike Duplicate (MS/MSD) for Oil & Grease method 1664B can only be analyzed when the client has submitted sufficient sample volume. An extra liter per MS/MSD is required to fulfill the method QC criteria. Please refer to Chain of Custody and QC Summary (MS/MSD) of the Laboratory Report to verify ample sample volume was submitted to fulfill the requirement.
pH	The method for pH does not stipulate a specific holding time other than to state that the samples should be analyzed as soon as possible. For aqueous samples the 40 CFR 136 specifies a holding time of 15 minutes from sampling to analysis. Therefore all aqueous pH samples not analyzed in the field are considered out of hold time at the time of sample receipt. All soil samples are analyzed as soon as possible after sample receipt.

Laboratory Control Sample (LCS): A known matrix spiked with compound(s) representative of the target analytes, which is used to document laboratory performance.

Matrix Duplicate: An intra-laboratory split sample which is used to document the precision of a method in a given sample matrix.

Matrix Spike: An aliquot of a sample spiked with a known concentration of target analyte(s). The spiking occurs prior to sample preparation and analysis. A matrix spike is used to document the bias of a method in a given sample matrix.

Method Blank: An analyte-free matrix to which all reagents are added in the same volumes or proportions as used in sample processing. The method blank should be carried through the complete sample preparation and analytical procedure. The method blank is used to document contamination resulting from the analytical process.

Method Detection Limit (MDL): The minimum concentration of a substance that can be measured and reported with 99% confidence that the analyte concentration is greater than zero and is determined from analysis of a sample in a given matrix type containing the analyte.

Reportable Detection Limit (RDL): The lowest concentration that can be reliably achieved within specified limits of precision and accuracy during routine laboratory operating conditions. For many analytes the RDL analyte concentration is selected as the lowest non-zero standard in the calibration curve. While the RDL is approximately 5 to 10 times the MDL, the RDL for each sample takes into account the sample volume/weight, extract/digestate volume, cleanup procedures and, if applicable, dry weight correction. Sample RDLs are highly matrix-dependent.

Surrogate: An organic compound which is similar to the target analyte(s) in chemical composition and behavior in the analytical process, but which is not normally found in environmental samples. These compounds are spiked into all blanks, standards, and samples prior to analysis. Percent recoveries are calculated for each surrogate.

Continuing Calibration Verification: The calibration relationship established during the initial calibration must be verified at periodic intervals. Concentrations, intervals, and criteria are method specific.

